## What is Claimed:

## 1. A compound selected from the formulas I or II:

$$R_1$$
 $R_2$ 
 $R_3$ 
 $R_4$ 
 $R_4$ 
 $R_5$ 
 $R_6$ 
 $C(CH_2)_{n-Y}$ 
 $R_6$ 
 $C(CH_2)_{n-Y}$ 
 $R_7$ 
 $R_8$ 
 $R_8$ 
 $R_9$ 
 $R_$ 

wherein:

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 $R_1$  is selected from H, OH or the  $C_1$ - $C_{12}$  esters (straight chain or branched) or  $C_1$ - $C_{12}$  (straight chain or branched or cyclic) alkyl ethers thereof, or halogens; or  $C_1$ - $C_4$  halogenated ethers including triflouromethyl ether and trichloromethyl ether.

 $R_2$ ,  $R_3$ ,  $R_4$ ,  $R_5$ , and  $R_6$  are independently selected from H, OH or the  $C_1$ - $C_{12}$  esters (straight chain or branched) or  $C_1$ - $C_{12}$  alkyl ethers (straight chain or branched or cyclic) thereof, halogens, or  $C_1$ - $C_4$  halogenated ethers including triflouromethyl ether and trichloromethyl ether, cyano,  $C_1$ - $C_6$  alkyl (straight chain or branched), or trifluoromethyl, with the proviso that, when  $R_1$  is H,  $R_2$  is not OH.

X is selected from H,  $C_1$ - $C_6$  alkyl, cyano, nitro, trifluoromethyl, halogen; n is 2 or 3;

Y is selected from:

a) the moiety:

wherein  $R_7$  and  $R_8$  are independently selected from the group of H,  $C_1$ - $C_6$  alkyl, or phenyl optionally substituted by CN,  $C_1$ - $C_6$  alkyl (straight chain or branched),  $C_1$ - $C_6$  alkoxy (straight chain or branched), halogen, -OH, -CF<sub>3</sub>, or -OCF<sub>3</sub>;

b) a five-membered saturated, unsaturated or partially unsaturated heterocycle containing up to two heteroatoms selected from the group consisting of -O-, -NH-, -N(C<sub>1</sub>C<sub>4</sub> alkyl)-, -N=, and -S(O)<sub>m</sub>-, wherein m is an integer of from 0-2, optionally substituted with 1-3 substituents independently selected from the group consisting of hydrogen, hydroxyl, halo, C<sub>1</sub>-C<sub>4</sub> alkyl, trihalomethyl, C<sub>1</sub>-C<sub>4</sub> alkoxy, trihalomethoxy, C<sub>1</sub>-C<sub>4</sub> acyloxy, C<sub>1</sub>-C<sub>4</sub> alkylthio, C<sub>1</sub>-C<sub>4</sub> alkylsulfinyl, C<sub>1</sub>-C<sub>4</sub> alkylsulfonyl, hydroxy (C<sub>1</sub>-C<sub>4</sub>)alkyl, -CO<sub>2</sub>H-, -CN-, -CONHR<sub>1</sub>-, -NH<sub>2</sub>-, C<sub>1</sub>-C<sub>4</sub> alkylamino, di(C<sub>1</sub>-C<sub>4</sub>)alkylamino, -NHSO<sub>2</sub>R<sub>1</sub>-, -NHCOR<sub>1</sub>-, -NO<sub>2</sub>, and phenyl optionally substituted with 1-3 (C<sub>1</sub>-C<sub>4</sub>)alkyl;

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c) a six-membered saturated, unsaturated or partially unsaturated heterocycle containing up to two heteroatoms selected from the group consisting of -O-, -NH-, -N( $C_1C_4$  alkyl)-, -N=, and -S(O)<sub>m</sub>-, wherein m is an integer of from 0-2, optionally substituted with 1-3 substituents independently selected from the group consisting of hydrogen, hydroxyl, halo,  $C_1$ - $C_4$  alkyl, trihalomethyl,  $C_1$ - $C_4$  alkoxy, trihalomethoxy,  $C_1$ - $C_4$  acyloxy,  $C_1$ - $C_4$  alkylthio,  $C_1$ - $C_4$  alkylsulfinyl,  $C_1$ - $C_4$  alkylsulfonyl, hydroxy ( $C_1$ - $C_4$ )alkyl, - $CO_2$ H-, -CN-, - $CONHR_1$ -, - $NH_2$ -,  $C_1$ - $C_4$  alkylamino, di( $C_1$ - $C_4$ )alkylamino, - $NHSO_2R_1$ -, - $NHCOR_1$ -, - $NO_2$ , and phenyl optionally substituted with 1-3 ( $C_1$ - $C_4$ )alkyl;

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d) a seven-membered saturated, unsaturated or partially unsaturated heterocycle containing up to two heteroatoms selected from the group consisting of -O-, -NH-, -N(C<sub>1</sub>C<sub>4</sub> alkyl)-, -N=, and -S(O)<sub>m</sub>-, wherein m is an integer of from 0-2, optionally substituted with 1-3 substituents independently selected from the group consisting of hydrogen, hydroxyl, halo, C<sub>1</sub>-C<sub>4</sub> alkyl, trihalomethyl, C<sub>1</sub>-C<sub>4</sub> alkoxy, trihalomethoxy, C<sub>1</sub>-C<sub>4</sub> acyloxy, C<sub>1</sub>-C<sub>4</sub> alkylthio, C<sub>1</sub>-C<sub>4</sub> alkylsulfinyl, C<sub>1</sub>-C<sub>4</sub> alkylsulfonyl, hydroxy (C<sub>1</sub>-C<sub>4</sub>)alkyl, -CO<sub>2</sub>H-, -CN-, -CONHR<sub>1</sub>-, -NH<sub>2</sub>-, C<sub>1</sub>-C<sub>4</sub> alkylamino, di(C<sub>1</sub>-C<sub>4</sub>)alkylamino, -NHSO<sub>2</sub>R<sub>1</sub>-, -NHCOR<sub>1</sub>-, -NO<sub>2</sub>, and phenyl optionally substituted with 1-3 (C<sub>1</sub>-C<sub>4</sub>)alkyl;; or

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e) a bicyclic heterocycle containing from 6-12 carbon atoms either bridged or fused and containing up to two heteroatoms selected from the group consisting of -O-, -NH-, -N( $C_1C_4$  alkyl)-, and -S(O)<sub>m</sub>-, wherein m is an integer of

from 0-2, optionally substituted with 1-3 substituents independently selected from the group consisting of hydrogen, hydroxyl, halo,  $C_1$ - $C_4$  alkyl, trihalomethyl,  $C_1$ - $C_4$  alkoxy, trihalomethoxy,  $C_1$ - $C_4$  acyloxy,  $C_1$ - $C_4$  alkylthio,  $C_1$ - $C_4$  alkylsulfinyl,  $C_1$ - $C_4$  alkylsulfonyl, hydroxy ( $C_1$ - $C_4$ )alkyl, - $CO_2$ H-, -CN-, - $CONHR_1$ -, - $NH_2$ -,  $C_1$ - $C_4$  alkylamino, di( $C_1$ - $C_4$ )alkylamino, - $NHSO_2R_1$ -, - $NHCOR_1$ -, - $NO_2$ , and phenyl optionally substituted with 1-3 ( $C_1$ - $C_4$ ) alkyl;

and the pharmaceutically acceptable salts thereof.

2. A compound of Claim 1 wherein:

 $R_1$  is selected from H, OH or the  $C_1$ - $C_4$  esters or alkyl ethers thereof, halogen;

 $R_2$ ,  $R_3$ ,  $R_4$ ,  $R_5$ , and  $R_6$  are independently selected from H, OH or the  $C_1$ - $C_4$  esters or alkyl ethers thereof, halogen, cyano,  $C_1$ - $C_6$  alkyl, or trifluoromethyl, with the proviso that, when  $R_1$  is H,  $R_2$  is not OH;

X is selected from H, C<sub>1</sub>-C<sub>6</sub> alkyl, cyano, nitro, triflouromethyl, halogen;

Y is the moiety

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R<sub>7</sub> and R<sub>8</sub> are selected independently from H, C<sub>1</sub>-C<sub>6</sub> alkyl, or combined by - (CH<sub>2</sub>)p-, wherein p is an integer of from 2 to 6, so as to form a ring, the ring being optionally substituted by up to three substituents selected from the group of hydrogen, hydroxyl, halo, C<sub>1</sub>-C<sub>4</sub> alkyl, trihalomethyl, C<sub>1</sub>-C<sub>4</sub> alkoxy, trihalomethoxy, C<sub>1</sub>-C<sub>4</sub> alkylthio, C<sub>1</sub>-C<sub>4</sub> alkylsulfinyl, C<sub>1</sub>-C<sub>4</sub> alkylsulfonyl, hydroxy (C<sub>1</sub>-C<sub>4</sub>)alkyl, -CO<sub>2</sub>H, -CN, -CONH(C<sub>1</sub>-C<sub>4</sub>), -NH<sub>2</sub>, C<sub>1</sub>-C<sub>4</sub> alkylamino, di(C<sub>1</sub>-C<sub>4</sub>)alkylamino, -NHSO<sub>2</sub>(C<sub>1</sub>-C<sub>4</sub>), -NHCO(C<sub>1</sub>-C<sub>4</sub>), and -NO<sub>2</sub>;

or a pharmaceutically acceptable salt thereof.

## 3. A compound of Claim 1 wherein:

R<sub>1</sub> is OH;

 $R_2$ ,  $R_3$ ,  $R_4$ ,  $R_5$ , and  $R_6$  are independently selected from H, OH or the  $C_1$ - $C_4$  esters or alkyl ethers thereof, halogen, cyano,  $C_1$ - $C_6$  alkyl, or trifluoromethyl, with the proviso that, when  $R_1$  is H,  $R_2$  is not OH;

X is selected from the group of Cl, NO<sub>2</sub>, CN, CF<sub>3</sub>, or CH<sub>3</sub>;

Y is the moiety

R<sub>7</sub> and R<sub>8</sub> are concatenated together as -(CH<sub>2</sub>)<sub>T</sub>, wherein r is an integer of from 4 to 6, to form a ring optionally substituted by up to three substituents selected from the group of hydrogen, hydroxyl, halo, C<sub>1</sub>-C<sub>4</sub> alkyl, trihalomethyl, C<sub>1</sub>-C<sub>4</sub> alkoxy trihalomethoxy, C<sub>1</sub> C<sub>4</sub> alkylthio, C<sub>1</sub>-C<sub>4</sub> alkylsulfinyl, C<sub>1</sub>-C<sub>4</sub> alkylsulfonyl, hydroxy (C<sub>1</sub>-C<sub>4</sub>)alkyl, -CO<sub>2</sub>H, -CN, -CONH(C<sub>1</sub>-C<sub>4</sub>), -NH<sub>2</sub>, C<sub>1</sub>-C<sub>4</sub> alkylamino, di(C<sub>1</sub>-C<sub>4</sub>)alkylamino, -NHSO<sub>2</sub>(C<sub>1</sub>-C<sub>4</sub>), -NHCO(C<sub>1</sub>-C<sub>4</sub>), and -NO<sub>2</sub>; or a pharmaceutically acceptable salt thereof.

- 4. A compound of Claim 1 which is 5-Benzyloxy-2-(4-ethoxy-phenyl)-3-methyl-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indole or a pharmaceutically acceptable salt thereof.
- 5. A compound of Claim 1 which is 5-Benzyloxy-2-phenyl-3-methyl-1-[4-(2-azepan-1-yl-ethoxy)-benzyl]-1H-indole or a pharmaceutically acceptable salt thereof.
- 6. A compound of Claim 1 which is 5-Benzyloxy-2-(4-benzyloxy-phenyl)3-methyl-1-[4-(2-azepan-1-yl-ethoxy)-benzyl]-1H-indole or a pharmaceutically acceptable salt thereof.
- 7. A compound of Claim 1 which is 5-Benzyloxy-2-(4-benzyloxy-phenyl)-3-methyl-1-[4-(2-diisopropylamino-1-yl-ethoxy)-benzyl]-1H-indole or a pharmaceutically acceptable salt thereof.
  - 8. A compound of Claim 1 which is 5-Benzyloxy-2-(4-benzyloxy-phenyl)-3-methyl-1-[4-(2-butyl-methylamino-1-ylethoxy)-benzyl]-1H-indole or a pharmaceutically acceptable salt thereof.

- 9. A compound of Claim 1 which is 5-Benzyloxy-2-(4-benzyloxy-phenyl)-3-methyl-1-{4-dimethylamino}-ethoxy}-benzyl}-1H-indole or a pharmaceutically acceptable salt thereof.
- 5 10. A compound of Claim 1 which is 5-Benzyloxy-2-(4-benzyloxy-phenyl)-3-methyl-1-{4-[2-(2-methyl-piperidin-1-yl)-ethoxy]-benzyl}-1H-indole or a pharmaceutically acceptable salt thereof.
- 11. A compound of Claum 1 which is 5-Benzyloxy-2-(4-benzyloxy-phenyl)3-methyl-1-{4-[2-(3-methyl-piperidin-1-yl)-ethoxy]-benzyl}-1H-indole or a
  pharmaceutically acceptable salt thereof.
- 12. A compound of Claim 1 which is 5-Benzyloxy-2-(4-benzyloxy-phenyl)-3-methyl-1-{4-[2-(4-methyl-piperidin-1-yl)-ethoxy]-benzyl}-1H-indole or a pharmaceutically acceptable salt thereof.
  - 13. A compound of Claim 1 which is 5-Benzyloxy-2-(4-benzyloxy-phenyl)-3-methyl-1{4-[2-((cis)-2,6-Dimethyl-piperidin-1-yl)-ethoxy]-benzyl}-1H-indole or a pharmaceutically acceptable salt thereof.
  - 14. A compound of Claim 1 which is 5-Benzyloxy-2-(4-benzyloxy-phenyl)-3-methyl-{4-[2-(1,3,3-trimethyl-6-aza-bicyclo[3.2.1]oct-6-yl)-ethoxy]-benzyl}-1H-indole or a pharmaceutically acceptable salt thereof.
- 25 15. A compound of Claim 1 which is (1S,4R)-5-Benzyloxy-2-(4-benzyloxy-phenyl)-3-methyl{4-[2-(2-Aza-bicyclo [2.2.1] hept-2-yl)-ethoxy]-benzyl}-1H-indole or a pharmaceutically acceptable salt thereof.

16. A compound of Claim 1 which is 5-Benzyloxy-2-(4-flouro-phenyl)-330 methyl-1-[4-(2-azepan-1-yl-ethoxy)-benzyl]-1H-indole or a pharmaceutically acceptable salt thereof.

- 17. A compound of Claim 1 which is 5-Benzyloxy-2-(4-flouro-phenyl)-3-methyl-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indole or a pharmaceutically acceptable salt thereof.
- 5 18. A compound of Claim 1 which is 5-Benzyloxy-2-(4-chloro-phenyl)-3-methyl-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indole or a pharmaceutically acceptable salt thereof.
- 19. A compound of Claim 1 which is 5-Benzyloxy-2-[3,4-methylenedioxy-10 phenyl]-3-methyl-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indole or a pharmaceutically acceptable salt thereof.
  - 20. A compound of Claim 1 which is 5-Benzyloxy-2-[4-isopropoxy-phenyl]-3-methyl-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indole or a pharmaceutically acceptable salt thereof.

- 21. A compound of Claim 1 which is 5-Benzyloxy-2-[4-methyl-phenyl]-3-methyl-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indole or a pharmaceutically acceptable salt thereof.
- 22. A compound of Claim 1 which is 1-[4-(2-Azepan-1-yl-ethoxy)-benzyl]-5-benzyloxy-2-(3-benzyloxy-phenyl)-3-methyl-1H-indole or a pharmaceutically acceptable salt thereof.
- 23. A compound of Claim 1 which is 5-Benzyloxy-2-(4-benzyloxy-3-fluoro-phenyl)-3-methyl-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indole or a pharmaceutically acceptable salt thereof.
- 24. A compound of Claim 1 which is 5-Benzyloxy-2-(4-benzyloxy-3-30 fluoro-phenyl)-3-methyl-1-[4-(2-azepan-1-yl-ethoxy)-benzyl]-1H-indole or a pharmaceutically acceptable salt thereof.

25. A compound of Claim 1 which is 5-Benzyloxy-2-(3-methoxy-phenyl-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-3-methyl-1H-indole or a pharmaceutically acceptable salt thereof.

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- 26. A compound of Claim 1 which is 5-Benzyloxy-3-methyl-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-2-(4-trifluoromethoxy-phenyl)-1H-indole or a pharmaceutically acceptable salt thereof.
- 27. A compound of Claim 1 which is (2-{4-[5-Benzyloxy-2-(4-benzyloxy-phenyl)-3-methyl-indol-1-ylmethyl]-phenoxy}-ethyl)-cyclohexyl-amine or a pharmaceutically acceptable salt thereof.
- 28. A compound of Claim 1 which is 5-Benzyloxy-2-(4-benzyloxy-phenyl)-15 3-methyl-1-{4-methylpiperazin-1-yl)-ethoxy]-benzyl}-1H-indole or a pharmaceutically acceptable salt thereof.
  - 29. A compound of Claim 1 which is 1-[4-(2-Azepan-1-yl-ethoxy)-benzyl]-5-benzyloxy-2-(3-methoxy-phenyl)-3-methyl-1H-indole or a pharmaceutically acceptable salt thereof.
  - 30. A compound of Claim 1 which is 4-{3-Methyl-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indole} (HCl).
- 25 31. A compound of Claim 1 which is 4-{3-Methyl-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indol-2-yl}-phenol hydrochloride (HCl).
  - 32. A compound of Claim 1 which is 3-Methyl-2-phenyl-1-[4-(2-piperidine-1-yl-ethoxy)-benzyl]-1H-indol-5-ol (HCl).

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33. A compound of Claim 1 which is 4-{5-Methoxy-3-methyl-1-{4-[2-(piperidin-1-yl)-ethoxy]-benzyl}-1H-indol-2-yl}-phenol or a pharmaceutically acceptable salt thereof.

34. A compound of Claim 1 which is 2-(4-methoxy-phenyl)-3-methyl-1-{4-[2-(piperidin-1-yl)-ethoxy]-benzyl}-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.

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- 35. A compound of Claim 1 which is 5-Methoxy-2-(4-methoxy-phenyl)-3-methyl-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indole (HCL).
- 36. A compound of Claim 1 which is 1-[4-(2-Azepan-1-yl-ethoxy)-benzyl]10 5-methoxy-2-(4-methoxy-phenyl)-3-methyl-1H-indole (HCL).
  - 37. A compound of Claim 1 which is 2-(4-Ethoxy-phenyl)-3-methyl-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.

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- 38. A compound of Claim 1 which is 1-[4-(2-Azepan-1-yl-ethoxy)-benzyl]-2-(4-ethoxy-phenyl)-3-methyl-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.
- 20 39. A compound of Claim 1 which is 4-{5-Fluoro-3-methyl-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indol-2-yl}-phenol (HCl).
  - 40. A compound of Claim 1 which is 1-[4-(2-Azepan-1-yl-ethoxy)-benzyl]-3-methyl-2-phenyl-1H-indol-5-ol (HCl).

- 41. A compound of Claim 1 which is 2-(4-Hydroxy-phenyl)-3-methyl-1-[4-(2-pyrollidin-1-yl-ethoxy)-benzyl]-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.
- 42. A compound of Claim 1 which is 1-[4-(2-Azepan-1-yl-ethoxy)-benzyl]-2-(4-hydroxy-phenyl)-3-methyl-1H-indol-5-ol (HCl).
  - 43. A compound of Claim 1 which is 1-[4-(2-Azepan-1-yl-ethoxy)-benzyl]-2-(4-hydroxy-phenyl)-3-methyl-1H-indol-5-ol Acetate Salt.

44. A compound of Claim 1 which is 1-[4-(2-Azocan-1-yl-ethoxy)-benzyl]-2-(4-hydroxy-phenyl)-3-methyl-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.

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- 45. A compound of Claim 1 which is 2-(4-Hydroxy-phenyl)-3-methyl-1-[4-(2-dimethyl-1-yl-ethoxy)-benzyl]-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.
- 10 46. A compound of Claim 1 which is 2-(4-Hydroxy-phenyl)-3-methyl-1-[4-(2-diethyl-1-yl-ethoxy)-benzyl]-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.
- 47. A compound of Claim 1 which is 1-[4-(2-Dipropylamino-ethoxy)-benzyl]-2-(4-hydroxy-phenyl)-3-methyl-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.
- 48. A compound of Claim 1 which is 1-[4-(2-Dibutylamino-ethoxy)-benzyl]-2-(4-hydroxy-phenyl)-3-methyl-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.
  - 49. A compound of Claim 1 which is 1-[4-(2-Diisopropylamino-ethoxy)-benzyl]-2-(4-hydroxy-phenyl)-3-methyl-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.

- 50. A compound of Claim 1 which is 1-{4-[2-(Butyl-methyl-amino)-ethoxy]-benzyl}-2-(4-hydroxy-phenyl)-3-methyl-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.
- 30 51. A compound of Claim 1 which is 2-(4-Hydroxy-phenyl)-3-methyl-1-{4-[2-(2-methyl-piperidin-1-yl)-ethoxy]-benzyl}-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.

- 52. A compound of Claim 1 which is 2-(4-Hydroxy-phenyl)-3-methyl-1-{4-[2-(3-methyl-piperdin-1-yl)-ethoxy]-benzyl}-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.
- 5 53. A compound of Claim 1 which is 2-(4-Hydroxy-phenyl)-3-methyl-1-{4-[2-(4-methyl-piperidin-1-yl)-ethoxy]-benzyl}-1H-indol-5-ol (HCl).
  - 54. A compound of Claim 1 which is 1-{4-[2-(3,3-Dimethyl-piperidin-1-yl)-ethoxy]-benzyl}-2-(4-hydroxy-phenyl)-3-methyl-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.

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- 55. A compound of Claim 1 which is 1-{4-[2-((cis)-2,6-Dimethyl-piperidin-1-yl)-ethoxy]-benzyl}-2-(4-hydroxy-phenyl)-3-methyl-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.
- 56. A compound of Claim 1 which is 2-(4-Hydroxy-phenyl)-1-{4-[2-(4-hydroxy-piperidin-1-yl)-ethoxy]-benzyl}-3-methyl-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.
- 20 57. A compound of Claim 1 which is (1S,4R)-1-{4-{2-(2-Aza-bicyclo [2.2.1] hept-2-yl)-ethoxy]-benzyl}-2-(4-hydroxy-phenyl)-3-methyl-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.
- 58. A compound of Claim 1 which is 2-(4-Hydroxy-phenyl)-3-methyl-125 {4-[2-(1,3,3-trimethyl-6-aza-bicyclo[3.2.1]oct-6-yl)-ethoxy]-benzyl}-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.
  - 59. A compound of Claim 1 which is 2-(4-Fluoro-phenyl)-3-methyl-1-[4-(2-piperidine-1-yl-ethoxy)-benzyl]-1H-indol-5-ol (HCl).
  - 60. A compound of Claim 1 which is 1-[4-(2-Azepan-1-yl-ethoxy)-benzyl]-2-(4-fluoro-phenyl)-3-methyl-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.

- 61. A compound of Claim 1 which is 2-(3-Methoxy-4-hydroxy-phenyl)-3-methyl-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indol-5-ol (HCl).
- 5 62. A compound of Claim 1 which is 2-Benzo[1,3]dioxol-5-yl-3-methyl-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indol-5-ol (HCL).
  - 63. A compound of Claim 1 which is 2-(4-Isopropoxy-phenyl)-3-methyl-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indol-5-ol (HCl).
  - 64. A compound of Claim 1 which is 1-[4-(2-Azepan-1-yl-ethoxy)-benzyl]-2-(4-isopropoxy-phenyl)-3-methyl-1H-indol-5-ol (HCl).
- 65. A compound of Claim 1 which is 2-(4-Cyclopenyloxy-phenyl)-3-methyl-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.

- 66. A compound of Claim 1 which is 3-Methyl-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-2-(4-trifluoromethyl-phenyl)-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.
  - 67. A compound of Claim 1 which is 3-Methyl-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-2-p-tolyl-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.
- 25 68. A compound of Claim 1 which is 2-(4-Chloro-phenyl)-3-methyl-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indol-5-ol (HCL).
- 69. A compound of Claim 1 which is 2-(2,4-Dimethoxy-phenyl)-3-methyl-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indol-5-ol or a pharmaceutically acceptable 30 salt thereof.

- 70. A compound of Claim 1 which is 2-(3-Hydroxy-phenyl)-3-methyl-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.
- 5 71. A compound of Claim 1 which is 1-[4-(2-Azepan-1-yl-ethoxy)-benzyl]-2-(3-hydroxy-phenyl)-3-methyl-1H-indole-5-ol or a pharmaceutically acceptable salt thereof.
- 72. A compound of Claim 1 which is 2-(3-Fluoro-4-hydroxy-phenyl)-3-10 methyl-1-[4-(2-pipendin-1-yl-ethoxy)-benzyl]-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.
  - 73. A compound of Claim 1 which is 2-(3-Fluoro-4-hydroxy-phenyl)-3-methyl-1-[4-(azepan-1-yl-ethoxy)-benzyl]-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.

- 74. A compound of Claim 1 which is 2-(3-Methoxy-phenyl)-3-methyl-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indole-5-ol or a pharmaceutically acceptable salt thereof.
- 75. A compound of Claim 1 which is 3-Methyl-1-[4-(2-piperidin-1-ylethoxy)-benzyl]-2-(4-trifluoromethoxy-phenyl)-1H-indole-5-ol or a pharmaceutically acceptable salt thereof.
- 25 76. A compound of Claim 1 which is 3-Chloro-2-(4-hydroxy-phenyl)-1-[4-(2-pyrrolidin-1-yl-ethoxy)-benzyl]-1H-indol-5-ol (HCl).
- 77. A compound of Claim 1 which is Removal of benzyl ethers to render 3-Chloro-2-(4-hydroxy-phenyl)-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indol-5-ol (HCl).
  - 78. A compound of Claim 1 which is 3-Chloro-2-(4-hydroxy-phenyl)-1-[4-(2-azepan-1-yl-ethoxy)-benzyl]-1H-indol-5-ol (HCl).

79. A compound of Claim 1 which is 3-Chloro-2-(4-hydroxy-2-methyl-phenyl)-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.

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- 80. A compound of Claim 1 which is 2-(4-Hydroxy-phenyl)-3-ethyl-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indol-5-ol (HCl).
- 81. A compound of Claim 1 which is 5-Hydroxy-2-(4-Hydroxy-phenyl)-1-10 [4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indole-3-carbonitrile (HCl).
  - 82. A compound of Claim 1 which is 1-[4-(2-Azepan-1-yl-ethoxy)-benzyl]-5-hydroxy-2-(4-hydroxy-phenyl)-1H-indole-3-cabonitrile (HCl).
- 15 83. A compound of Claim 1 which is 5-Benzyloxy-2-(4-benzyloxy-phenyl)-3-chloro-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indole or a pharmaceutically acceptable salt thereof.
- 84. A compound of Claim 1 which is 5-Benzyloxy-2-(4-benzyloxy-phenyl)3-chloro-1-[4-(2-azepan-1-yl-ethoxy)-benzyl]-1H-indole or a pharmaceutically acceptable salt thereof.
- 85. A compound of Claim 1 which is 5-Benzyloxy-2-(2-methyl-4-benzyloxy-phenyl)-3-chloro-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indole or a pharmaceutically acceptable salt thereof.
  - 86. A compound of Claim 1 which is 5-Benzyloxy-2-(4-benzyloxy-phenyl)-3-ethyl-1-[4-(2-piperidin -1-yl-ethoxy)-benzyl]-1H-indole or a pharmaceutically acceptable salt thereof.
  - 87. A compound of Claim 1 which is 5-Benzyloxy-2-(4-benzyloxy-phenyl)-3-cyano-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indole or a pharmaceutically acceptable salt thereof.

88. A compound of Claim 1 which is 5-Benzyloxy-2-(4-benzyloxy-phenyl)-3-cyano-1-[4-(2-azepan-1-yl-ethoxy)-benzyl]-1H-indole or a pharmaceutically acceptable salt thereof.

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- 89. A compound of Claim 1 which is Di-propionate of 1-[4-(2-Azepan-1-ylethoxy)-benzyl]-2-(4-hydroxy-phenyl)-3-methyl-1H-indol-5-ol (HCl).
- 90. A compound of Claim 1 which is Di-pivalate of 1-[4-(2-Azepan-1-yl-10 ethoxy)-benzyl]-2-(4-hydroxy-phenyl)-3-methyl-1H-indol-5-ol (HCl).
  - 91. A compound of Claim 1 which is 5-Benzyloxy-2-(4-benzyloxy-phenyl)-1-[4-(3-piperidin-1-yl-propoxy)-benzyl]-3-methyl-1H-indole or a pharmaceutically acceptable salt thereof.

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- 92. A compound of Claim 1 which is 2-(4-Hydroxy-phenyl)-3-methyl-1-{4-[3-(piperidin-1-yl)-propoxy]-benzyl}-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.
- 93. A compound of Claim 1 which is 2-(4-Hydroxy-phenyl)-1-[3-methoxy-4-(2-piperidin-1-yl-ethoxy)-benzyl]-3-methyl-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.
- 94. A compound of Claim 1 which is 2-(4-Hydroxy-phenyl)-1-[3-methoxy-4-(2-azepan-1-yl-ethoxy)-benzyl]-3-methyl-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.
  - 95. A compound of Claim 1 which is 5-Benzyloxy-2-(4-benzyloxy-phenyl)-3-methyl-1-[3-Methoxy-4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indole or a pharmaceutically acceptable salt thereof.

- 96. A compound of Claim 1 which is 5-Benzyloxy-2-(4-benzyloxy-phenyl)-3-methyl-1-[2-Methoxy-4-(2-azepan-1-yl-ethoxy)-benzyl]-1H-indole or a pharmaceutically acceptable salt thereof.
- 5 97. A compound of Claim 1 which is Di-pivalate ester of 2-(4-Hydroxy-phenyl)-3-methyl-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.
- 98. A compound of Claim 1 which is 5-Benzyloxy-2-(4-benzyloxy-phenyt)3-methyl-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indole or a pharmaceutically acceptable salt thereof.
- 99. A compound of Claim 1 which is 5-Benzyloxy-2-(3-benzyloxy-phenyl)-3-methyl-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indole or a pharmaceutically acceptable salt thereof.
  - 100. A compound of Claim 1 which is 2-(4-Hydroxy-phenyl)-3-methyl-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indol-5-ol or a pharmaceutically acceptable salt thereof.

101. A compound of Claim 1 which is 2-(4-Hydroxy-phenyl)-3-methyl-1-[4-(2-piperidin-1-yl-ethoxy)-benzyl]-1H-indol-5-ol methiodide.

- 102. A compound of Claim 1 which is 2-(4-Hydroxy-phenyl)-3-methyl-1-[4-(2-dimethyl-1-yl-ethoxy)-benzyl]-1H-indol-5-ol methiodide.
  - 103. A pharmaceutical composition comprising a compound of Claim 1, or a pharmaceutically acceptable salt thereof, and a pharmaceutically acceptable carrier or excipient.

104. A method of treating or preventing bone loss in a mammal, the method comprising administering to a mammal in need thereof an effective amount of a compound of Claim 1, or a pharmaceutically acceptable salt thereof.

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- 105. A method of treating or preventing disease states or syndromes which are caused or associated with an estrogen deficiency in a mammal, the method comprising administering to a mammal in need thereof an effective amount of a compound of Claim 1, or a pharmaceutically acceptable salt thereof.
- 106. A method of treating or preventing cardiovascular disease in a mammal, the method comprising administering to a mammal in need thereof an effective amount of a compound of Claim 1, or a pharmaccutically acceptable salt thereof.

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107. A method of treating or preventing disease in a mammal which result from proliferation or abnormal development, actions or growth of endometrial or endometrial-like tissue, the method comprising administering to a mammal in need thereof an effective amount of a compound of Claim 1, or a pharmaceutically acceptable salt thereof.